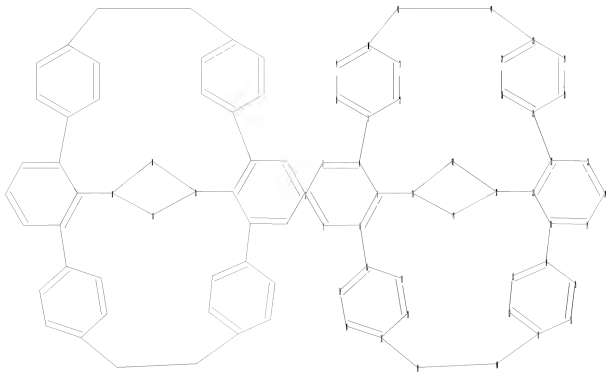

ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23
24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39 40 41 42 43 44
ring bonds :
1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-37 6-16 7-8 7-12 8-9 9-10 10-11 10-43
11-12 13-14 13-18 13-39 14-15 15-16 16-17 17-18 19-20 19-24 20-21 20-34
21-22 21-38 22-23 22-25 23-24 25-26 25-30 26-27 27-28 28-29 28-44 29-30
31-32 31-36 31-40 32-33 33-34 34-35 35-36 37-41 37-42 38-41 38-42 39-40
43-44
exact/norm bonds :
4-7 5-37 6-16 10-43 13-39 20-34 21-38 22-25 28-44 31-40 37-41 37-42
38-41 38-42 39-40 43-44
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-12 8-9 9-10 10-11 11-12 13-14 13-18
14-15 15-16 16-17 17-18 19-20 19-24 20-21 21-22 22-23 23-24 25-26 25-30
26-27 27-28 28-29 29-30 31-32 31-36 32-33 33-34 34-35 35-36

Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom
20:Atom 21:Atom 22:Atom 23:Atom 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom
29:Atom 30:Atom 31:Atom 32:Atom 33:Atom 34:Atom 35:Atom 36:Atom 37:Atom
38:Atom 39:Atom 40:Atom 41:Atom 42:Atom 43:Atom 44:Atom

L1 STRUCTURE UPLOADED

```
=> d his

     (FILE 'HOME' ENTERED AT 12:42:48 ON 30 MAR 2008)

     FILE 'REGISTRY' ENTERED AT 12:43:04 ON 30 MAR 2008
L1             STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1             STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> s l1
SAMPLE SEARCH INITIATED 12:43:33 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -     4688 TO ITERATE

 42.7% PROCESSED     2000 ITERATIONS                                 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:        89654 TO    97866
PROJECTED ANSWERS:               0 TO        0

L2            0 SEA SSS SAM L1

=> s l1 full
FULL SEARCH INITIATED 12:43:40 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -    93556 TO ITERATE

100.0% PROCESSED    93556 ITERATIONS                                0 ANSWERS
SEARCH TIME: 00.00.01

L3            0 SEA SSS FUL L1

=> fil caplus
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                       ENTRY    SESSION
FULL ESTIMATED COST                                   179.28     179.49

FILE 'CAPLUS' ENTERED AT 12:44:58 ON 30 MAR 2008
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FILE COVERS 1907 - 30 Mar 2008  VOL 148 ISS 14
FILE LAST UPDATED: 28 Mar 2008  (20080328/ED)

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They are available for your review at:

http://www.cas.org/infopolicy.html

=> s cyclophane#
L4         4837 CYCLOPHANE#

=> s l4 and diamines
          29064 DIAMINES
L5           25 L4 AND DIAMINES

=> s l5 and complex
       1400406 COMPLEX
L6            5 L5 AND COMPLEX

=> d 1-5 bib aabs
'AABS' IS NOT A VALID FORMAT FOR FILE 'CAPLUS'

The following are valid formats:

ABS --------- GI and AB
ALL --------- BIB, AB, IND, RE
APPS -------- AI, PRAI
BIB --------- AN, plus Bibliographic Data and PI table (default)
CAN --------- List of CA abstract numbers without answer numbers
CBIB -------- AN, plus Compressed Bibliographic Data
CLASS ------- IPC, NCL, ECLA, FTERM
DALL -------- ALL, delimited (end of each field identified)
DMAX -------- MAX, delimited for post-processing
FAM --------- AN, PI and PRAI in table, plus Patent Family data
FBIB -------- AN, BIB, plus Patent FAM
IND --------- Indexing data
IPC --------- International Patent Classifications
MAX --------- ALL, plus Patent FAM, RE
PATS -------- PI, SO
SAM --------- CC, SX, TI, ST, IT
SCAN -------- CC, SX, TI, ST, IT (random display, no answer numbers;
              SCAN must be entered on the same line as the DISPLAY,
              e.g., D SCAN or DISPLAY SCAN)
STD --------- BIB, CLASS

IABS -------- ABS, indented with text labels
IALL -------- ALL, indented with text labels
IBIB -------- BIB, indented with text labels
IMAX -------- MAX, indented with text labels
ISTD -------- STD, indented with text labels

OBIB -------- AN, plus Bibliographic Data (original)
OIBIB ------- OBIB, indented with text labels

SBIB -------- BIB, no citations
SIBIB ------- IBIB, no citations

HIT --------- Fields containing hit terms
HITIND ------ IC, ICA, ICI, NCL, CC and index field (ST and IT)
              containing hit terms
HITRN ------- HIT RN and its text modification
HITSTR ------ HIT RN, its text modification, its CA index name, and
```

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             its structure diagram
HITSEQ ------ HIT RN, its text modification, its CA index name, its
             structure diagram, plus NTE and SEQ fields
FHITSTR ----- First HIT RN, its text modification, its CA index name, and
             its structure diagram
FHITSEQ ----- First HIT RN, its text modification, its CA index name, its
             structure diagram, plus NTE and SEQ fields
KWIC -------- Hit term plus 20 words on either side
OCC --------- Number of occurrence of hit term and field in which it occurs

To display a particular field or fields, enter the display field
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an arrow prompt (=>).  Examples of formats include: TI; TI,AU; BIB,ST;
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specification.

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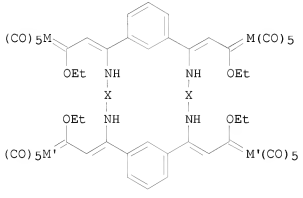
to view a specified Accession Number.
ENTER DISPLAY FORMAT (BIB):end

=> d 1-5 bib abs

L6   ANSWER 1 OF 5  CAPLUS  COPYRIGHT 2008 ACS on STN
AN   2006:1079231  CAPLUS
DN   146:27919
TI   Structure-Activity Relationships for Cytotoxic Ruthenium(II) Arene
     Complexes Containing N,N-, N,O-, and O,O-Chelating Ligands
AU   Habtemariam, Abraha; Melchart, Michael; Fernandez, Rafael; Parsons, Simon;
     Oswald, Iain D. H.; Parkin, Andrew; Fabbiani, Francesca P. A.; Davidson,
     James E.; Dawson, Alice; Aird, Rhona E.; Jodrell, Duncan I.; Sadler, Peter
     J.
CS   School of Chemistry, University of Edinburgh, Edinburgh, EH9 3JJ, UK
SO   Journal of Medicinal Chemistry (2006), 49(23), 6858-6868
     CODEN: JMCMAR; ISSN: 0022-2623
PB   American Chemical Society
DT   Journal
LA   English
OS   CASREACT 146:27919
AB   Ruthenium arene complexes containing bidentate diamine, amino acid and
     diketonate chelate ligands were prepared by a variety of appropriate
     procedures and examined for cytostatic activity against human cancer cells.
     Organometallic Ru(II) complexes [(η6-arene)Ru(XY)Cl]Z, where XY is an
     N,N- (diamine), N,O- (e.g., amino acidate), or O,O- (e.g.,
     β-diketonate) chelating ligand, the arene ranges from benzene derivs.
     to fused polycyclic hydrocarbons, and Z is usually PF6, were prepared by
     direct or reduction-assisted complexation of arenes, substitution of
     cycloalkadiene or arene ligands with subsequent complexation of bidentate
     XY-ligands.  The x-ray structures of 13 complexes are reported.  All have
     the characteristic "piano-stool" geometry.  The structure-activity
     relationships was evaluated for cytotoxicity of the prepared complexes
     against human cancer cells.  The complexes most active toward A2780 human
     ovarian cancer cells contained XY = ethylenediamine (en) and extended
     polycyclic arenes.  Complexes with polar substituents on the arene or XY =
     bipyridyl derivs. exhibited reduced activity.  The activity of the
     O,O-chelated complexes depended strongly on the substituents and on the
     arene.  For arene = p-cymene, XY = amino acidate complexes were inactive.
     Complexes were not cross-resistant with cisplatin, and cross-resistance to
     Adriamycin was circumvented by replacing XY = en with 1,2-
     phenylenediamine.  Some complexes were also active against colon,
     pancreatic, and lung cancer cells.
```

RE.CNT 88 THERE ARE 88 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 2 OF 5 CAPLUS COPYRIGHT 2008 ACS on STN
AN 2003:213301 CAPLUS
DN 138:368989
TI Synthesis and Electrochemical Properties of Novel Tetrametallic Macrocyclic Fischer Carbene Complexes
AU Fernandez, Israel; Mancheno, Maria Jose; Gomez-Gallego, Mar; Sierra, Miguel A.
CS Departamento de Quimica Organica, Facultad de Quimica, Universidad Complutense de Madrid, Madrid, 28040, Spain
SO Organic Letters (2003), 5(8), 1237-1240
CODEN: ORLEF7; ISSN: 1523-7060
PB American Chemical Society
DT Journal
LA English
OS CASREACT 138:368989
GI




I

AB Metallomacrocyclic compds. [I; wherein M, M', independently = Cr, W; X = -C6H4-, -(C6H4)2-] can be easily prepared by 1,4-addition of diamines to α,β-unsatd. Fischer bis-carbene templates. This method allows the preparation of a new family of homo- and heterotetrametallic compds. having macrocyclic cyclophanic structures.
RE.CNT 45 THERE ARE 45 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 3 OF 5 CAPLUS COPYRIGHT 2008 ACS on STN
AN 2001:680382 CAPLUS
DN 136:6093
TI Synthesis of Cyclophanic Chromium(0) Bis(carbene) Complexes
AU Fernandez, Israel; Sierra, Miguel A.; Mancheno, Maria Jose; Gomez-Gallego, Mar; Ricart, Susagna
CS Departamento de Quimica Organica Facultad de Quimica, Universidad Complutense, Madrid, 28040, Spain
SO Organometallics (2001), 20(21), 4304-4306
CODEN: ORGND7; ISSN: 0276-7333
PB American Chemical Society
DT Journal
LA English
OS CASREACT 136:6093

AB Double Michael addition of 1,4- and 1,3-xylylenediamines and 1,5-diaminopentane to the bimetallic α,β-unsatd. alkoxychromium(0) carbene complex m-C6H4[C.tplbond.CC(OEt):Cr(CO)5]2 4 produces in good to excellent yields the corresponding cyclophane bimetallic complexes as single isomers. In contrast, the reaction with 1,4-diaminobutane produces a different cyclophane complex, derived from a 1,2- + 1,4-addition process.

RE.CNT 33 THERE ARE 33 CITED REFERENCES AVAILABLE FOR THIS RECORD
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L6 ANSWER 4 OF 5 CAPLUS COPYRIGHT 2008 ACS on STN
AN 1988:131780 CAPLUS
DN 108:131780
TI Polyaza-macrocycles of cyclophane type: synthesis, structure of a chloroform inclusion complex and anion binding
AU Jazwinski, Jaroslaw; Lehn, Jean Marie; Meric, Robert; Vigneron, Jean Pierre; Cesario, Michele; Guilhem, J.; Pascard, Claudine
CS Inst. Bel, Univ. Louis Pasteur, Strasbourg, Fed. Rep. Ger.
SO Tetrahedron Letters (1987), 28(30), 3489-92
CODEN: TELEAY; ISSN: 0040-4039
DT Journal
LA English
OS CASREACT 108:131780
GI

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

AB Cyclocondensation of 4-HCOC6H4CH2C6H4CHO-4 with diamines, e.g., H2NCH2CH2NH2, (H2NCH2CH2)2X (X = O, NSO2C6H4Me-4) gave 65-85% I and II, resp. I gave a CHCl3 inclusion complex, the crystal structure of which was determined Reduction of I by LiAlH4 gave tetramine III, which binds dicarboxylate substrates.

L6 ANSWER 5 OF 5 CAPLUS COPYRIGHT 2008 ACS on STN
AN 1984:434686 CAPLUS
DN 101:34686
OREF 101:5369a,5372a
TI Cyclophane porphyrins and their metal complexes. Biomimetic study on receptor site of hemoprotein oxygen binding
AU Ogoshi, Hisanobu; Sugimoto, Hiroshi; Miyake, Masao; Yoshida, Zenichi
CS Fac. Eng., Kyoto Univ., Kyoto, 606, Japan
SO Tetrahedron (1984), 40(3), 579-92
CODEN: TETRAB; ISSN: 0040-4020
DT Journal
LA English
AB A new sym. porphyrin, 7,8,17,18-tetraethyl-3,13-dimethylporphyrin-2,12-dipropionic acid, and its derivs. were synthesized by the a,c-biladiene route. Condensation of the dipropionic acid with diamine, [H2N(CH2)nNH2, n = 6, 7, 8, 9, 10, 12, and 14], afforded the corresponding cyclophane porphyrins. The bridged groups were characterized by the 1H NMR spectra of their Zn complexes. The spin state of the Fe3+ complexes of the cyclophane porphyrins was investigated by changing the size of the bridged chain or size of axial ligand. The cyclophane-porphyrinato(III) perchlorate complexes in a mixture of MeOH and CHCl3 with 4-benzylpyridine provide a model for methemoproteins. Steric constraint between and axial ligand and the bridge group, [CH2CH2CONH(CH2)nNHCOCH2CH2] at the bridged face dets. the ratio of the

penta- and hexa-coordinated ferric complexes. The rate of O-binding to the Co2+ cyclophane porphyrins is markedly dependent on the size of the bridge chain. Removal of a solvent mol. or 6th axial ligand from the near proximity of the Co2+ complex apparently increases the rate of O binding.

=>

---Logging off of STN---

=>
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| DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) | SINCE FILE ENTRY | TOTAL SESSION |
| CA SUBSCRIBER PRICE | -4.00 | -4.00 |

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